



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 5, 2023 – 03:58 AM EDT

PDB ID : 6V6K
Title : EGFR(T790M/V948R) in complex with LN2057
Authors : Heppner, D.E.; Eck, M.J.
Deposited on : 2019-12-05
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : **FAILED**
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : **FAILED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

There are no overall percentile quality scores available for this entry.

MolProbity and EDS failed to run properly - the sequence quality summary graphics cannot be shown.

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	QQJ	F	1202	-	X	-	-
3	QQJ	H	1202	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 20750 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Epidermal growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	D	314	Total 2521	C 1614	N 427	O 461	S 19	0	0	0
1	A	300	Total 2420	C 1553	N 408	O 440	S 19	0	1	0
1	B	301	Total 2424	C 1554	N 411	O 440	S 19	0	0	0
1	C	302	Total 2440	C 1564	N 413	O 444	S 19	0	1	0
1	E	303	Total 2441	C 1565	N 413	O 444	S 19	0	1	0
1	F	306	Total 2453	C 1572	N 416	O 446	S 19	0	0	0
1	G	314	Total 2530	C 1619	N 428	O 464	S 19	0	1	0
1	H	314	Total 2530	C 1619	N 428	O 464	S 19	0	1	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	790	MET	THR	engineered mutation	UNP P00533
D	948	ARG	VAL	engineered mutation	UNP P00533
A	790	MET	THR	engineered mutation	UNP P00533
A	948	ARG	VAL	engineered mutation	UNP P00533
B	790	MET	THR	engineered mutation	UNP P00533
B	948	ARG	VAL	engineered mutation	UNP P00533
C	790	MET	THR	engineered mutation	UNP P00533
C	948	ARG	VAL	engineered mutation	UNP P00533
E	790	MET	THR	engineered mutation	UNP P00533
E	948	ARG	VAL	engineered mutation	UNP P00533
F	790	MET	THR	engineered mutation	UNP P00533
F	948	ARG	VAL	engineered mutation	UNP P00533
G	790	MET	THR	engineered mutation	UNP P00533

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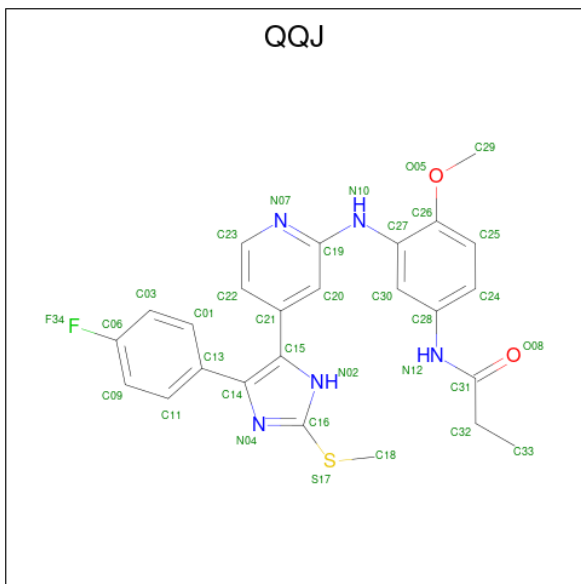
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Chain	Residue	Modelled	Actual	Comment	Reference
G	948	ARG	VAL	engineered mutation	UNP P00533
H	790	MET	THR	engineered mutation	UNP P00533
H	948	ARG	VAL	engineered mutation	UNP P00533

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total Cl 1 1	0	0
2	A	1	Total Cl 1 1	0	0
2	B	1	Total Cl 1 1	0	0
2	C	1	Total Cl 1 1	0	0
2	E	1	Total Cl 1 1	0	0
2	F	1	Total Cl 1 1	0	0
2	G	1	Total Cl 1 1	0	0
2	H	1	Total Cl 1 1	0	0

- Molecule 3 is N-[3-({4-[4-(4-fluorophenyl)-2-(methylsulfanyl)-1H-imidazol-5-yl]pyridin-2-yl}amino)-4-methoxyphenyl]propanamide (three-letter code: QQJ) (formula: C₂₅H₂₄FN₅O₂S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
3	D	1	Total	C	F	N	O	S	0	0
			34	25	1	5	2	1		
3	A	1	Total	C	F	N	O	S	0	0
			34	25	1	5	2	1		
3	B	1	Total	C	F	N	O	S	0	0
			34	25	1	5	2	1		
3	C	1	Total	C	F	N	O	S	0	0
			34	25	1	5	2	1		
3	E	1	Total	C	F	N	O	S	0	0
			34	25	1	5	2	1		
3	F	1	Total	C	F	N	O	S	0	0
			34	25	1	5	2	1		
3	G	1	Total	C	F	N	O	S	0	0
			34	25	1	5	2	1		
3	H	1	Total	C	F	N	O	S	0	0
			34	25	1	5	2	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	87	Total	O	0	0
			87	87		
4	A	81	Total	O	0	0
			81	81		
4	B	80	Total	O	0	0
			80	80		
4	C	103	Total	O	0	0
			103	103		
4	E	78	Total	O	0	0
			78	78		
4	F	85	Total	O	0	0
			85	85		
4	G	86	Total	O	0	0
			86	86		
4	H	111	Total	O	0	0
			111	111		

MolProbity and EDS failed to run properly - this section is therefore empty.

3 Data and refinement statistics i

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.60Å 102.45Å 174.04Å 90.00° 101.25° 90.00°	Depositor
Resolution (Å)	85.35 – 2.20	Depositor
% Data completeness (in resolution range)	93.7 (85.35-2.20)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.20Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.214 , 0.239	Depositor
Wilson B-factor (Å ²)	27.7	Xtrriage
Anisotropy	0.717	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.005 for h,-k,-h-l	Xtrriage
Total number of atoms	20750	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 66.40 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.0839e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

4 Model quality [i](#)

4.1 Standard geometry [i](#)

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4.2 Too-close contacts [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3.2 Protein sidechains [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3.3 RNA [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

4.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

4.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 8 are monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	QQJ	G	1202	1	35,37,37	5.92	28 (80%)	42,51,51	1.62	9 (21%)
3	QQJ	D	1202	1	35,37,37	5.82	30 (85%)	42,51,51	1.61	10 (23%)
3	QQJ	E	1202	1	35,37,37	4.75	27 (77%)	42,51,51	2.65	12 (28%)
3	QQJ	F	1202	1	35,37,37	5.74	28 (80%)	42,51,51	1.71	12 (28%)
3	QQJ	C	1202	1	35,37,37	5.80	30 (85%)	42,51,51	1.60	11 (26%)
3	QQJ	B	1202	1	35,37,37	4.79	29 (82%)	42,51,51	2.00	9 (21%)
3	QQJ	H	1202	1	35,37,37	5.97	30 (85%)	42,51,51	2.07	10 (23%)
3	QQJ	A	1202	1	35,37,37	5.71	29 (82%)	42,51,51	1.73	8 (19%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	QQJ	G	1202	1	-	4/20/22/22	0/4/4/4
3	QQJ	D	1202	1	-	4/20/22/22	0/4/4/4
3	QQJ	E	1202	1	-	4/20/22/22	0/4/4/4
3	QQJ	F	1202	1	-	6/20/22/22	0/4/4/4
3	QQJ	C	1202	1	-	2/20/22/22	0/4/4/4
3	QQJ	B	1202	1	-	1/20/22/22	0/4/4/4
3	QQJ	H	1202	1	-	6/20/22/22	0/4/4/4
3	QQJ	A	1202	1	-	2/20/22/22	0/4/4/4

All (231) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1202	QQJ	C30-C28	9.96	1.55	1.39
3	C	1202	QQJ	C30-C28	9.91	1.55	1.39
3	A	1202	QQJ	C30-C28	9.61	1.55	1.39
3	G	1202	QQJ	C30-C28	9.54	1.55	1.39
3	C	1202	QQJ	C30-C27	9.51	1.54	1.39
3	F	1202	QQJ	C30-C27	9.45	1.54	1.39
3	H	1202	QQJ	C30-C28	9.41	1.54	1.39
3	G	1202	QQJ	C24-C28	9.37	1.55	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	1202	QQJ	C30-C27	9.35	1.54	1.39
3	A	1202	QQJ	C24-C28	9.27	1.54	1.39
3	F	1202	QQJ	C30-C28	9.27	1.54	1.39
3	D	1202	QQJ	C30-C27	9.20	1.53	1.39
3	F	1202	QQJ	C25-C24	9.16	1.55	1.38
3	D	1202	QQJ	C24-C28	9.09	1.54	1.39
3	A	1202	QQJ	C30-C27	9.08	1.53	1.39
3	F	1202	QQJ	C24-C28	9.03	1.54	1.39
3	H	1202	QQJ	C30-C27	9.02	1.53	1.39
3	A	1202	QQJ	C25-C24	9.00	1.55	1.38
3	H	1202	QQJ	C24-C28	8.99	1.54	1.39
3	G	1202	QQJ	C25-C24	8.94	1.55	1.38
3	E	1202	QQJ	C30-C28	8.91	1.54	1.39
3	C	1202	QQJ	C25-C24	8.85	1.54	1.38
3	H	1202	QQJ	C25-C24	8.84	1.54	1.38
3	C	1202	QQJ	C24-C28	8.77	1.54	1.39
3	B	1202	QQJ	C25-C24	8.49	1.54	1.38
3	D	1202	QQJ	C25-C24	8.48	1.54	1.38
3	G	1202	QQJ	C16-S17	8.46	1.82	1.75
3	H	1202	QQJ	C22-C23	8.19	1.55	1.38
3	H	1202	QQJ	C16-S17	8.11	1.82	1.75
3	H	1202	QQJ	C03-C06	8.04	1.52	1.37
3	G	1202	QQJ	C03-C06	8.01	1.52	1.37
3	C	1202	QQJ	C03-C06	7.99	1.52	1.37
3	H	1202	QQJ	C03-C01	7.99	1.53	1.38
3	F	1202	QQJ	C03-C06	7.98	1.52	1.37
3	E	1202	QQJ	C25-C24	7.92	1.53	1.38
3	D	1202	QQJ	C03-C06	7.90	1.52	1.37
3	A	1202	QQJ	C03-C06	7.88	1.52	1.37
3	C	1202	QQJ	C03-C01	7.87	1.53	1.38
3	A	1202	QQJ	C03-C01	7.84	1.53	1.38
3	G	1202	QQJ	C09-C06	7.83	1.52	1.37
3	A	1202	QQJ	C22-C23	7.79	1.54	1.38
3	G	1202	QQJ	C22-C23	7.76	1.54	1.38
3	D	1202	QQJ	C22-C23	7.75	1.54	1.38
3	G	1202	QQJ	C09-C11	7.73	1.52	1.38
3	H	1202	QQJ	C09-C11	7.71	1.52	1.38
3	G	1202	QQJ	C03-C01	7.70	1.52	1.38
3	D	1202	QQJ	C09-C11	7.68	1.52	1.38
3	F	1202	QQJ	C09-C11	7.68	1.52	1.38
3	C	1202	QQJ	C09-C11	7.67	1.52	1.38
3	B	1202	QQJ	C30-C28	7.67	1.52	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1202	QQJ	C03-C01	7.63	1.52	1.38
3	E	1202	QQJ	C30-C27	7.63	1.51	1.39
3	F	1202	QQJ	C03-C01	7.62	1.52	1.38
3	B	1202	QQJ	C30-C27	7.58	1.51	1.39
3	H	1202	QQJ	C23-N07	7.58	1.50	1.34
3	D	1202	QQJ	C09-C06	7.58	1.52	1.37
3	C	1202	QQJ	C22-C23	7.58	1.53	1.38
3	F	1202	QQJ	C22-C23	7.58	1.53	1.38
3	D	1202	QQJ	C16-S17	7.50	1.82	1.75
3	H	1202	QQJ	C19-N07	7.50	1.48	1.34
3	H	1202	QQJ	C09-C06	7.48	1.51	1.37
3	F	1202	QQJ	C09-C06	7.46	1.51	1.37
3	C	1202	QQJ	C09-C06	7.43	1.51	1.37
3	D	1202	QQJ	C23-N07	7.40	1.50	1.34
3	A	1202	QQJ	C09-C11	7.40	1.52	1.38
3	C	1202	QQJ	C23-N07	7.37	1.50	1.34
3	E	1202	QQJ	C24-C28	7.35	1.51	1.39
3	A	1202	QQJ	C16-S17	7.29	1.81	1.75
3	F	1202	QQJ	C25-C26	7.26	1.54	1.39
3	A	1202	QQJ	C09-C06	7.26	1.51	1.37
3	G	1202	QQJ	C23-N07	7.23	1.49	1.34
3	F	1202	QQJ	C23-N07	7.21	1.49	1.34
3	F	1202	QQJ	C16-S17	7.21	1.81	1.75
3	A	1202	QQJ	C23-N07	7.21	1.49	1.34
3	H	1202	QQJ	C25-C26	7.17	1.54	1.39
3	B	1202	QQJ	C24-C28	7.16	1.51	1.39
3	G	1202	QQJ	C19-N07	7.16	1.47	1.34
3	G	1202	QQJ	C25-C26	7.08	1.54	1.39
3	A	1202	QQJ	C25-C26	7.06	1.54	1.39
3	C	1202	QQJ	C25-C26	7.03	1.54	1.39
3	D	1202	QQJ	C25-C26	7.02	1.54	1.39
3	A	1202	QQJ	C19-N07	7.00	1.47	1.34
3	C	1202	QQJ	C19-N07	6.97	1.47	1.34
3	F	1202	QQJ	C19-N07	6.93	1.47	1.34
3	D	1202	QQJ	C19-N07	6.91	1.47	1.34
3	B	1202	QQJ	C22-C23	6.69	1.52	1.38
3	G	1202	QQJ	C11-C13	6.51	1.53	1.39
3	B	1202	QQJ	C33-C32	-6.49	1.22	1.51
3	H	1202	QQJ	C11-C13	6.48	1.53	1.39
3	D	1202	QQJ	C11-C13	6.47	1.53	1.39
3	H	1202	QQJ	C20-C19	6.41	1.55	1.39
3	H	1202	QQJ	C20-C21	6.30	1.50	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	1202	QQJ	C20-C19	6.28	1.55	1.39
3	C	1202	QQJ	C16-S17	6.28	1.81	1.75
3	A	1202	QQJ	C11-C13	6.22	1.52	1.39
3	C	1202	QQJ	C11-C13	6.19	1.52	1.39
3	C	1202	QQJ	C33-C32	-6.18	1.23	1.51
3	D	1202	QQJ	C20-C19	6.15	1.55	1.39
3	B	1202	QQJ	O08-C31	-6.14	1.10	1.23
3	G	1202	QQJ	C20-C21	6.06	1.50	1.39
3	E	1202	QQJ	C22-C23	6.06	1.50	1.38
3	D	1202	QQJ	C20-C21	6.04	1.50	1.39
3	E	1202	QQJ	C09-C11	6.04	1.49	1.38
3	G	1202	QQJ	C01-C13	6.00	1.52	1.39
3	E	1202	QQJ	C16-S17	5.99	1.80	1.75
3	F	1202	QQJ	C11-C13	5.95	1.52	1.39
3	E	1202	QQJ	C03-C01	5.92	1.49	1.38
3	H	1202	QQJ	C27-C26	5.90	1.53	1.40
3	A	1202	QQJ	C20-C19	5.88	1.54	1.39
3	D	1202	QQJ	C01-C13	5.84	1.51	1.39
3	F	1202	QQJ	C01-C13	5.82	1.51	1.39
3	B	1202	QQJ	C03-C01	5.79	1.49	1.38
3	C	1202	QQJ	C20-C19	5.79	1.54	1.39
3	C	1202	QQJ	C01-C13	5.79	1.51	1.39
3	H	1202	QQJ	C01-C13	5.79	1.51	1.39
3	H	1202	QQJ	C21-C15	5.79	1.55	1.49
3	F	1202	QQJ	C20-C21	5.78	1.49	1.39
3	F	1202	QQJ	C20-C19	5.78	1.54	1.39
3	E	1202	QQJ	C03-C06	5.74	1.48	1.37
3	B	1202	QQJ	C25-C26	5.68	1.51	1.39
3	A	1202	QQJ	C01-C13	5.67	1.51	1.39
3	E	1202	QQJ	C25-C26	5.66	1.51	1.39
3	G	1202	QQJ	C31-N12	5.65	1.48	1.35
3	B	1202	QQJ	C03-C06	5.64	1.48	1.37
3	B	1202	QQJ	C15-N02	-5.64	1.23	1.37
3	F	1202	QQJ	C31-N12	5.63	1.48	1.35
3	E	1202	QQJ	C23-N07	5.53	1.46	1.34
3	C	1202	QQJ	C20-C21	5.52	1.49	1.39
3	E	1202	QQJ	C20-C19	5.52	1.53	1.39
3	B	1202	QQJ	C09-C06	5.51	1.47	1.37
3	A	1202	QQJ	C27-C26	5.50	1.52	1.40
3	D	1202	QQJ	C27-C26	5.50	1.52	1.40
3	H	1202	QQJ	C31-N12	5.50	1.47	1.35
3	F	1202	QQJ	C27-C26	5.41	1.52	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1202	QQJ	C20-C21	5.41	1.49	1.39
3	C	1202	QQJ	C27-C26	5.38	1.51	1.40
3	D	1202	QQJ	C31-N12	5.37	1.47	1.35
3	G	1202	QQJ	C27-C26	5.36	1.51	1.40
3	B	1202	QQJ	C19-N07	5.27	1.44	1.34
3	C	1202	QQJ	C31-N12	5.23	1.47	1.35
3	A	1202	QQJ	C31-N12	5.22	1.47	1.35
3	E	1202	QQJ	C09-C06	5.22	1.47	1.37
3	B	1202	QQJ	C20-C19	5.21	1.52	1.39
3	B	1202	QQJ	C09-C11	5.19	1.48	1.38
3	H	1202	QQJ	C19-N10	5.13	1.47	1.38
3	E	1202	QQJ	C19-N07	5.13	1.44	1.34
3	E	1202	QQJ	O08-C31	-5.09	1.12	1.23
3	D	1202	QQJ	C19-N10	5.08	1.47	1.38
3	G	1202	QQJ	C19-N10	4.90	1.47	1.38
3	B	1202	QQJ	C11-C13	4.89	1.49	1.39
3	C	1202	QQJ	C19-N10	4.88	1.47	1.38
3	H	1202	QQJ	C22-C21	4.79	1.49	1.39
3	G	1202	QQJ	C21-C15	4.71	1.54	1.49
3	E	1202	QQJ	C20-C21	4.68	1.47	1.39
3	E	1202	QQJ	C15-N02	-4.67	1.26	1.37
3	B	1202	QQJ	C23-N07	4.60	1.44	1.34
3	F	1202	QQJ	C19-N10	4.59	1.46	1.38
3	G	1202	QQJ	C22-C21	4.56	1.49	1.39
3	F	1202	QQJ	C22-C21	4.55	1.49	1.39
3	D	1202	QQJ	C22-C21	4.53	1.49	1.39
3	D	1202	QQJ	C13-C14	4.48	1.54	1.49
3	G	1202	QQJ	C13-C14	4.47	1.54	1.49
3	B	1202	QQJ	C14-N04	-4.41	1.26	1.37
3	B	1202	QQJ	C16-S17	4.39	1.79	1.75
3	E	1202	QQJ	C27-C26	4.38	1.49	1.40
3	B	1202	QQJ	C15-C14	-4.38	1.32	1.44
3	A	1202	QQJ	C19-N10	4.36	1.46	1.38
3	A	1202	QQJ	C22-C21	4.34	1.48	1.39
3	B	1202	QQJ	C01-C13	4.33	1.48	1.39
3	D	1202	QQJ	C21-C15	4.29	1.54	1.49
3	C	1202	QQJ	C22-C21	4.28	1.48	1.39
3	H	1202	QQJ	C13-C14	4.27	1.54	1.49
3	E	1202	QQJ	C11-C13	4.17	1.48	1.39
3	E	1202	QQJ	C01-C13	4.15	1.48	1.39
3	B	1202	QQJ	C27-C26	4.04	1.49	1.40
3	E	1202	QQJ	C15-C14	-4.00	1.33	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1202	QQJ	C21-C15	4.00	1.53	1.49
3	E	1202	QQJ	C14-N04	-3.98	1.27	1.37
3	A	1202	QQJ	C13-C14	3.98	1.53	1.49
3	F	1202	QQJ	C21-C15	3.79	1.53	1.49
3	B	1202	QQJ	C22-C21	3.74	1.47	1.39
3	A	1202	QQJ	C21-C15	3.65	1.53	1.49
3	F	1202	QQJ	C28-N12	3.63	1.49	1.41
3	E	1202	QQJ	C13-C14	3.60	1.53	1.49
3	D	1202	QQJ	C27-N10	3.27	1.49	1.39
3	F	1202	QQJ	C13-C14	3.23	1.52	1.49
3	G	1202	QQJ	C28-N12	3.13	1.48	1.41
3	C	1202	QQJ	C28-N12	3.08	1.47	1.41
3	B	1202	QQJ	C31-N12	3.06	1.42	1.35
3	H	1202	QQJ	C27-N10	3.04	1.48	1.39
3	C	1202	QQJ	C13-C14	2.97	1.52	1.49
3	A	1202	QQJ	C28-N12	2.94	1.47	1.41
3	C	1202	QQJ	C15-C14	-2.94	1.36	1.44
3	A	1202	QQJ	C15-C14	-2.94	1.36	1.44
3	C	1202	QQJ	C27-N10	2.94	1.48	1.39
3	E	1202	QQJ	C22-C21	2.92	1.45	1.39
3	G	1202	QQJ	C27-N10	2.91	1.48	1.39
3	E	1202	QQJ	C31-N12	2.88	1.42	1.35
3	D	1202	QQJ	C28-N12	2.87	1.47	1.41
3	F	1202	QQJ	C15-C14	-2.86	1.36	1.44
3	G	1202	QQJ	O05-C26	2.86	1.41	1.37
3	H	1202	QQJ	C28-N12	2.84	1.47	1.41
3	A	1202	QQJ	C27-N10	2.83	1.47	1.39
3	F	1202	QQJ	C27-N10	2.82	1.47	1.39
3	H	1202	QQJ	C33-C32	-2.76	1.39	1.51
3	G	1202	QQJ	C15-C14	-2.74	1.37	1.44
3	H	1202	QQJ	C15-C14	-2.72	1.37	1.44
3	E	1202	QQJ	C21-C15	2.69	1.52	1.49
3	D	1202	QQJ	C15-C14	-2.59	1.37	1.44
3	B	1202	QQJ	C20-C21	2.56	1.44	1.39
3	D	1202	QQJ	C15-N02	-2.54	1.31	1.37
3	C	1202	QQJ	O05-C26	2.52	1.41	1.37
3	B	1202	QQJ	C13-C14	2.47	1.51	1.49
3	G	1202	QQJ	C15-N02	-2.45	1.31	1.37
3	B	1202	QQJ	F34-C06	-2.39	1.30	1.36
3	F	1202	QQJ	O05-C26	2.35	1.40	1.37
3	F	1202	QQJ	C15-N02	-2.33	1.31	1.37
3	B	1202	QQJ	C32-C31	2.32	1.58	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	1202	QQJ	O08-C31	-2.29	1.18	1.23
3	A	1202	QQJ	O05-C26	2.29	1.40	1.37
3	C	1202	QQJ	C15-N02	-2.26	1.31	1.37
3	H	1202	QQJ	C15-N02	-2.26	1.31	1.37
3	H	1202	QQJ	O05-C26	2.23	1.40	1.37
3	A	1202	QQJ	C15-N02	-2.21	1.31	1.37
3	B	1202	QQJ	O05-C29	-2.19	1.36	1.42
3	A	1202	QQJ	C14-N04	-2.18	1.32	1.37
3	E	1202	QQJ	F34-C06	-2.18	1.31	1.36
3	D	1202	QQJ	C14-N04	-2.06	1.32	1.37
3	D	1202	QQJ	O05-C26	2.03	1.40	1.37
3	D	1202	QQJ	O08-C31	-2.01	1.19	1.23
3	C	1202	QQJ	C14-N04	-2.00	1.32	1.37

All (81) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	1202	QQJ	C18-S17-C16	12.83	111.85	102.27
3	H	1202	QQJ	C18-S17-C16	8.17	108.36	102.27
3	B	1202	QQJ	C18-S17-C16	7.59	107.93	102.27
3	A	1202	QQJ	C18-S17-C16	6.16	106.86	102.27
3	H	1202	QQJ	C28-N12-C31	-5.11	118.56	127.50
3	E	1202	QQJ	C28-N12-C31	-5.01	118.73	127.50
3	F	1202	QQJ	C18-S17-C16	5.00	106.00	102.27
3	B	1202	QQJ	O05-C26-C27	4.38	120.18	114.80
3	H	1202	QQJ	O05-C26-C27	4.05	119.77	114.80
3	B	1202	QQJ	O08-C31-N12	-3.97	116.38	123.63
3	B	1202	QQJ	C33-C32-C31	3.87	125.58	113.30
3	E	1202	QQJ	O05-C26-C27	3.80	119.46	114.80
3	G	1202	QQJ	C18-S17-C16	3.76	105.08	102.27
3	A	1202	QQJ	C22-C23-N07	-3.55	119.55	123.96
3	D	1202	QQJ	C20-C19-N07	-3.49	117.89	122.75
3	D	1202	QQJ	C23-N07-C19	3.30	121.90	117.22
3	E	1202	QQJ	C15-C14-N04	-3.28	104.92	113.76
3	B	1202	QQJ	C22-C23-N07	-3.23	119.94	123.96
3	G	1202	QQJ	C22-C23-N07	-3.22	119.96	123.96
3	C	1202	QQJ	O05-C26-C27	3.21	118.74	114.80
3	C	1202	QQJ	C22-C23-N07	-3.10	120.10	123.96
3	F	1202	QQJ	C28-N12-C31	-3.07	122.13	127.50
3	G	1202	QQJ	C23-N07-C19	3.04	121.53	117.22
3	G	1202	QQJ	O05-C26-C27	3.01	118.49	114.80
3	H	1202	QQJ	C22-C23-N07	-2.99	120.25	123.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	1202	QQJ	C28-N12-C31	-2.97	122.30	127.50
3	E	1202	QQJ	C22-C23-N07	-2.95	120.29	123.96
3	F	1202	QQJ	C29-O05-C26	-2.84	113.25	117.53
3	C	1202	QQJ	C29-O05-C26	-2.83	113.26	117.53
3	D	1202	QQJ	C20-C21-C15	-2.83	114.71	120.15
3	G	1202	QQJ	C20-C19-N07	-2.78	118.88	122.75
3	F	1202	QQJ	C20-C19-N07	-2.78	118.88	122.75
3	C	1202	QQJ	C23-N07-C19	2.76	121.14	117.22
3	E	1202	QQJ	C22-C21-C15	-2.76	116.24	120.61
3	D	1202	QQJ	C22-C21-C20	2.75	122.05	118.16
3	H	1202	QQJ	C29-O05-C26	-2.70	113.46	117.53
3	D	1202	QQJ	C15-C14-N04	-2.65	106.62	113.76
3	H	1202	QQJ	O05-C26-C25	-2.63	119.86	124.37
3	F	1202	QQJ	C23-N07-C19	2.63	120.95	117.22
3	A	1202	QQJ	C23-N07-C19	2.62	120.94	117.22
3	C	1202	QQJ	C20-C19-N07	-2.62	119.10	122.75
3	A	1202	QQJ	O05-C26-C27	2.61	118.00	114.80
3	F	1202	QQJ	C22-C21-C20	2.59	121.82	118.16
3	C	1202	QQJ	C22-C21-C20	2.57	121.80	118.16
3	A	1202	QQJ	C22-C21-C20	2.56	121.79	118.16
3	F	1202	QQJ	O05-C26-C27	2.55	117.93	114.80
3	E	1202	QQJ	C23-N07-C19	2.53	120.81	117.22
3	D	1202	QQJ	C22-C23-N07	-2.52	120.82	123.96
3	H	1202	QQJ	C20-C19-N07	-2.52	119.23	122.75
3	E	1202	QQJ	C09-C06-C03	-2.47	119.55	122.83
3	H	1202	QQJ	C23-N07-C19	2.47	120.72	117.22
3	C	1202	QQJ	C33-C32-C31	-2.46	105.49	113.30
3	G	1202	QQJ	C15-C14-N04	-2.44	107.16	113.76
3	G	1202	QQJ	C33-C32-C31	-2.42	105.60	113.30
3	A	1202	QQJ	C15-C14-N04	-2.42	107.24	113.76
3	D	1202	QQJ	O05-C26-C25	-2.38	120.29	124.37
3	E	1202	QQJ	F34-C06-C03	2.37	122.57	118.54
3	B	1202	QQJ	C15-C14-N04	-2.35	107.42	113.76
3	F	1202	QQJ	C22-C23-N07	-2.31	121.09	123.96
3	D	1202	QQJ	C18-S17-C16	2.27	103.96	102.27
3	C	1202	QQJ	C11-C13-C01	2.24	122.06	117.59
3	F	1202	QQJ	C15-C14-N04	-2.24	107.71	113.76
3	D	1202	QQJ	O05-C26-C27	2.24	117.54	114.80
3	B	1202	QQJ	C11-C09-C06	2.24	120.67	118.36
3	D	1202	QQJ	C28-N12-C31	-2.23	123.59	127.50
3	A	1202	QQJ	C11-C13-C01	2.21	121.99	117.59
3	H	1202	QQJ	C14-C15-N02	-2.17	107.90	113.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	1202	QQJ	C15-C14-N04	-2.17	107.91	113.76
3	A	1202	QQJ	C29-O05-C26	-2.16	114.27	117.53
3	F	1202	QQJ	C32-C31-N12	2.15	120.51	114.75
3	G	1202	QQJ	C22-C21-C20	2.13	121.18	118.16
3	F	1202	QQJ	C11-C13-C01	2.13	121.84	117.59
3	F	1202	QQJ	C20-C21-C15	-2.11	116.09	120.15
3	B	1202	QQJ	O05-C26-C25	-2.10	120.76	124.37
3	E	1202	QQJ	C22-C21-C20	2.10	121.13	118.16
3	E	1202	QQJ	C01-C13-C14	2.09	123.92	120.61
3	C	1202	QQJ	C15-C14-N04	-2.09	108.13	113.76
3	B	1202	QQJ	C14-C15-N02	-2.08	108.14	113.76
3	E	1202	QQJ	O05-C26-C25	-2.08	120.81	124.37
3	C	1202	QQJ	O05-C26-C25	-2.08	120.81	124.37
3	C	1202	QQJ	C20-C21-C15	-2.04	116.23	120.15

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1202	QQJ	C01-C13-C14-N04
3	H	1202	QQJ	C27-C26-O05-C29
3	F	1202	QQJ	C27-C26-O05-C29
3	H	1202	QQJ	C25-C26-O05-C29
3	D	1202	QQJ	C25-C26-O05-C29
3	F	1202	QQJ	C25-C26-O05-C29
3	D	1202	QQJ	C27-C26-O05-C29
3	G	1202	QQJ	C01-C13-C14-N04
3	E	1202	QQJ	N12-C31-C32-C33
3	A	1202	QQJ	C11-C13-C14-N04
3	G	1202	QQJ	C11-C13-C14-N04
3	H	1202	QQJ	C01-C13-C14-N04
3	H	1202	QQJ	C11-C13-C14-N04
3	F	1202	QQJ	N12-C31-C32-C33
3	C	1202	QQJ	C01-C13-C14-N04
3	C	1202	QQJ	C11-C13-C14-N04
3	F	1202	QQJ	C01-C13-C14-N04
3	F	1202	QQJ	C11-C13-C14-N04
3	E	1202	QQJ	O08-C31-C32-C33
3	F	1202	QQJ	O08-C31-C32-C33
3	H	1202	QQJ	O08-C31-C32-C33
3	D	1202	QQJ	C11-C13-C14-N04
3	G	1202	QQJ	N12-C31-C32-C33

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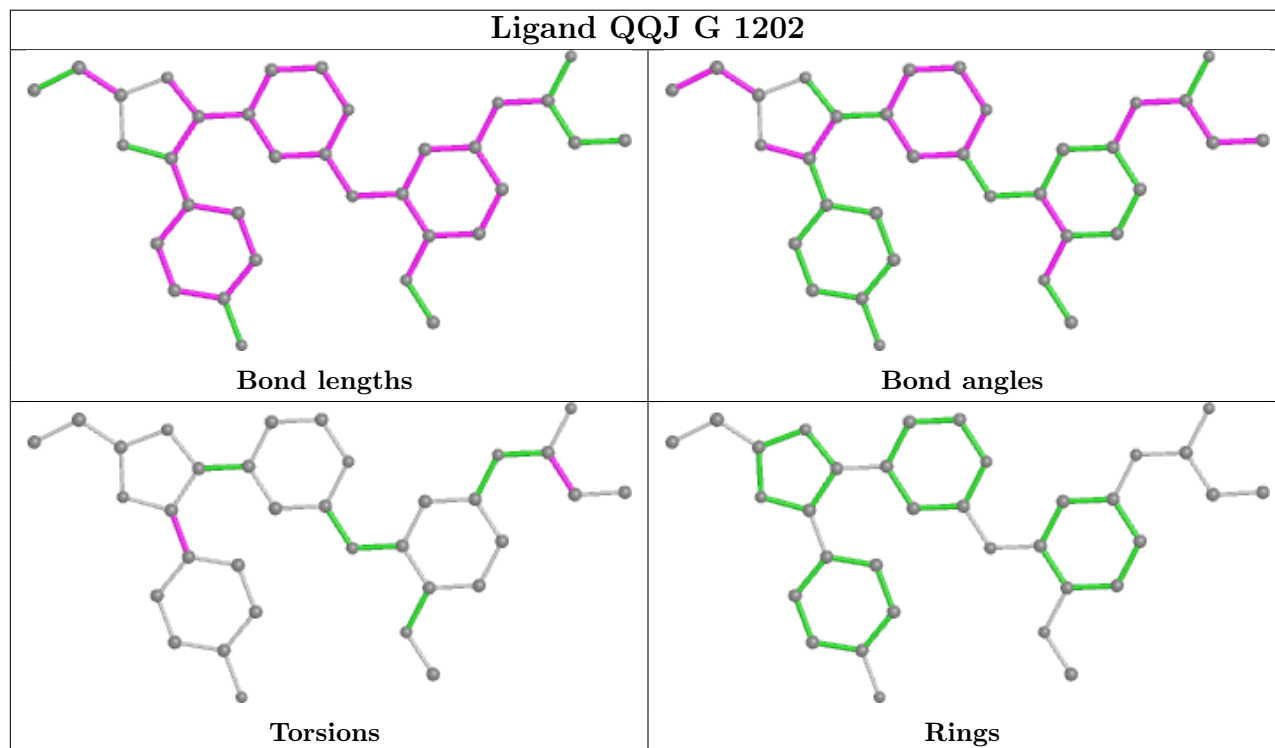
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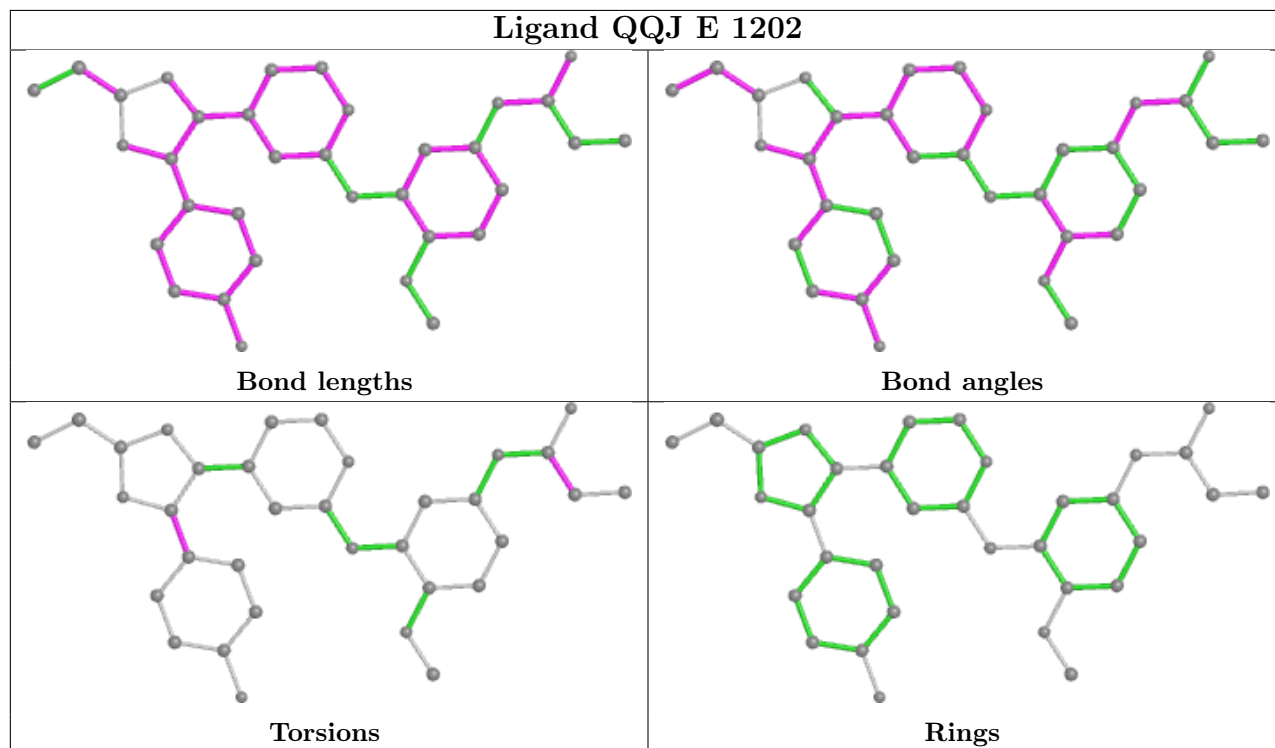
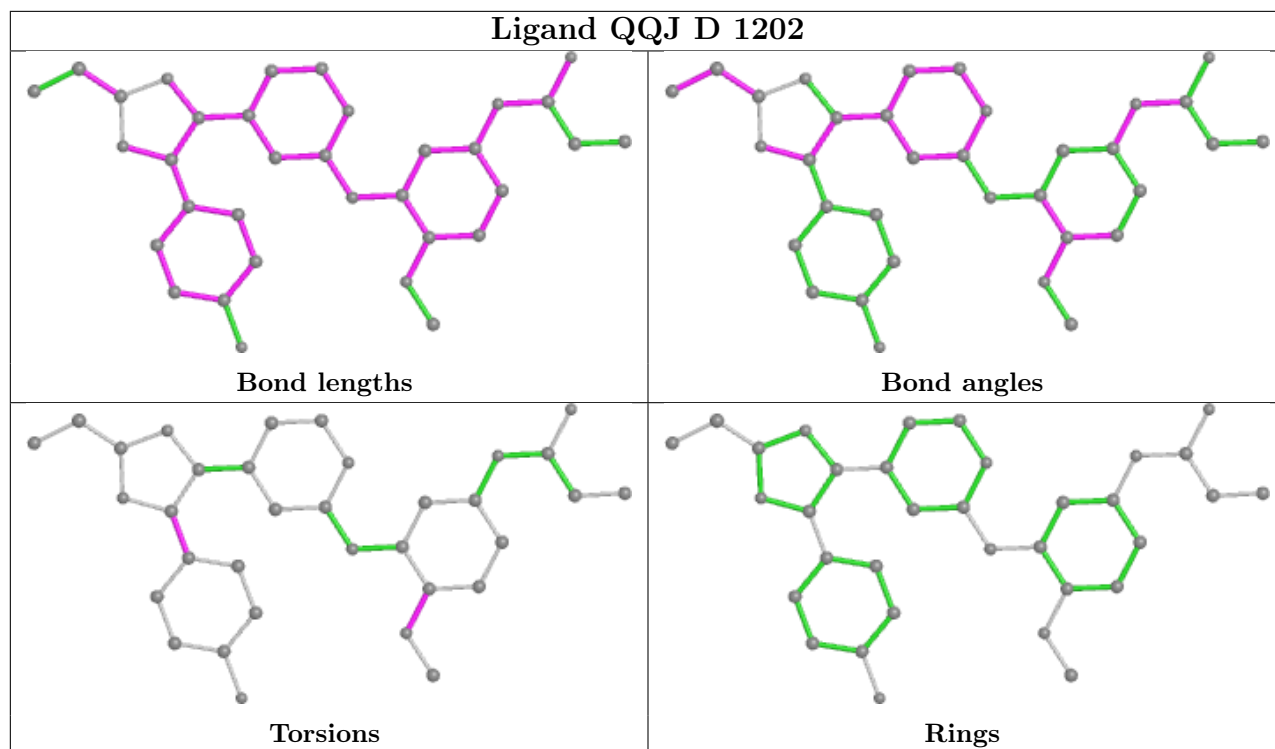
Mol	Chain	Res	Type	Atoms
3	G	1202	QQJ	O08-C31-C32-C33
3	D	1202	QQJ	C01-C13-C14-N04
3	B	1202	QQJ	C01-C13-C14-N04
3	E	1202	QQJ	C01-C13-C14-N04
3	E	1202	QQJ	C11-C13-C14-N04
3	H	1202	QQJ	N12-C31-C32-C33

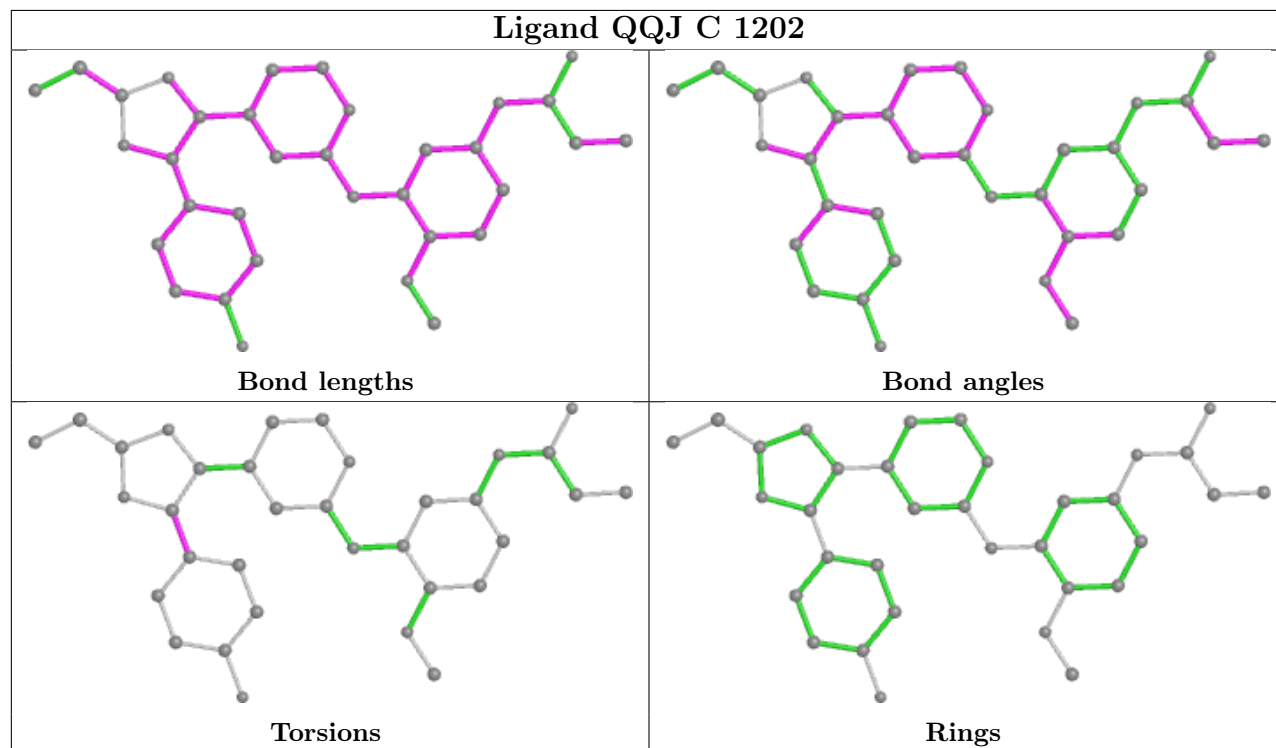
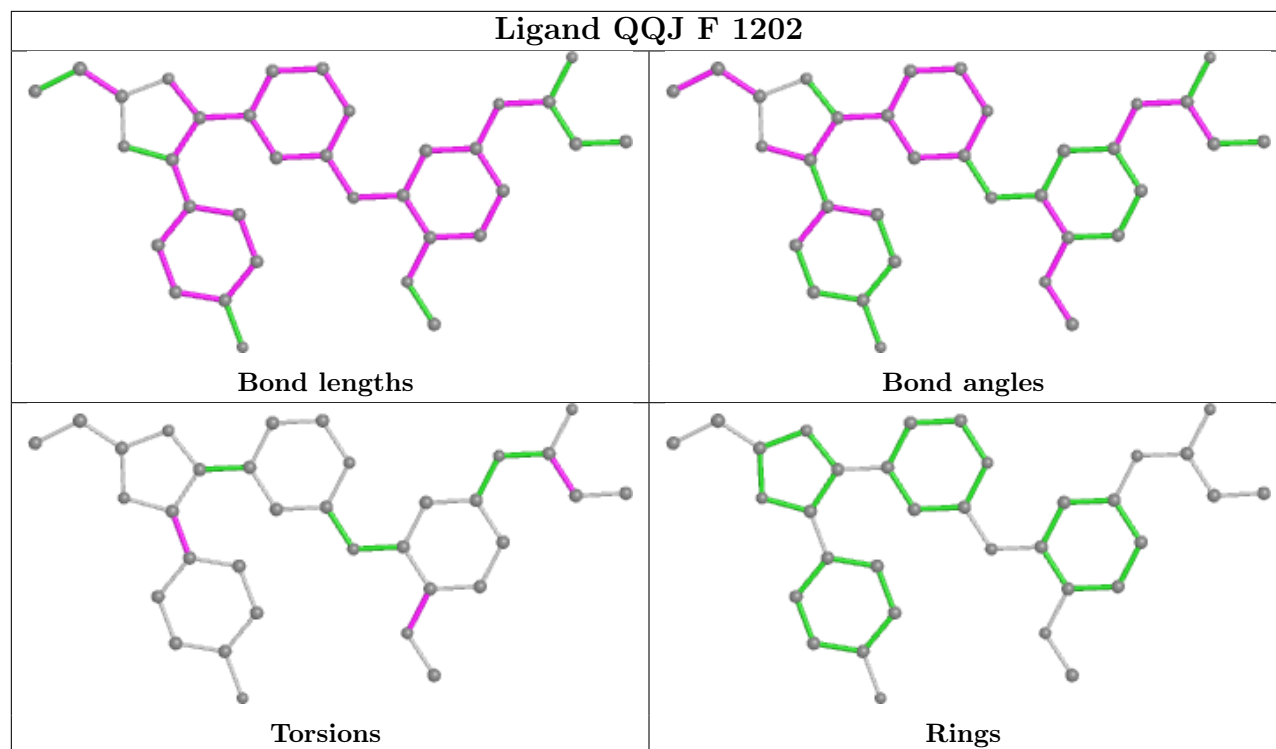
There are no ring outliers.

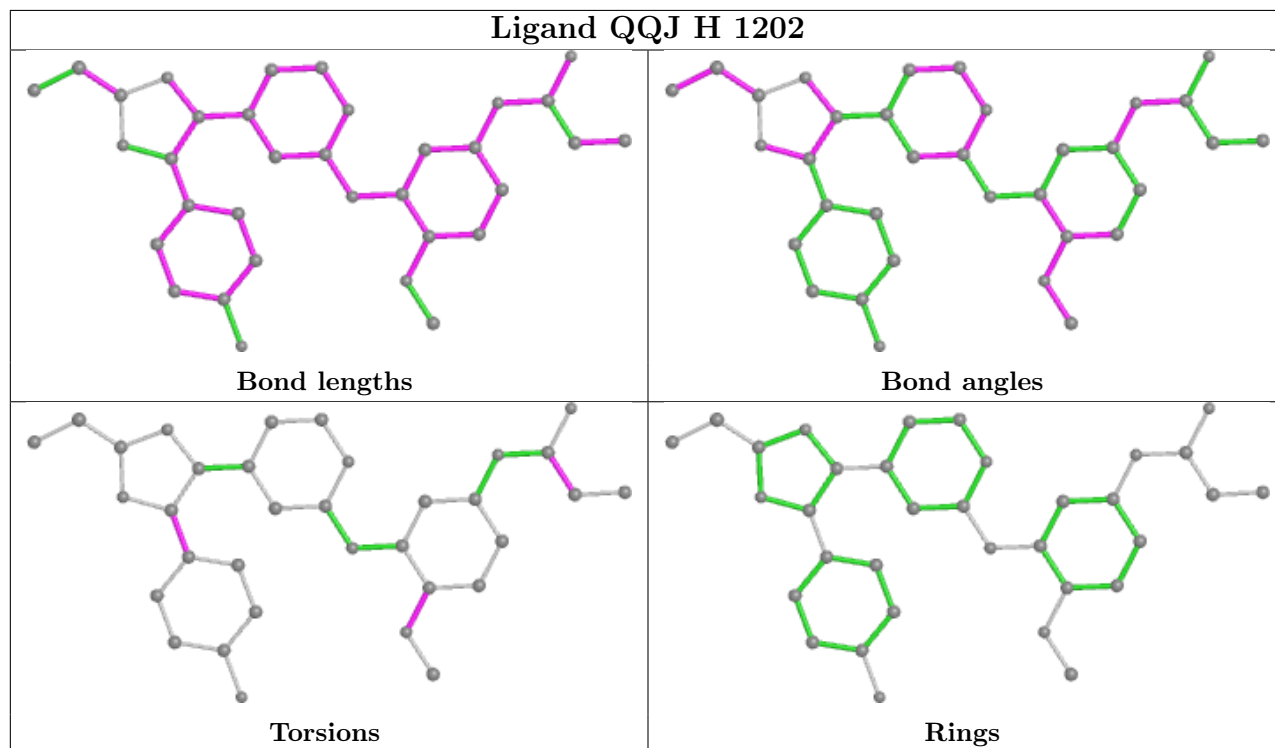
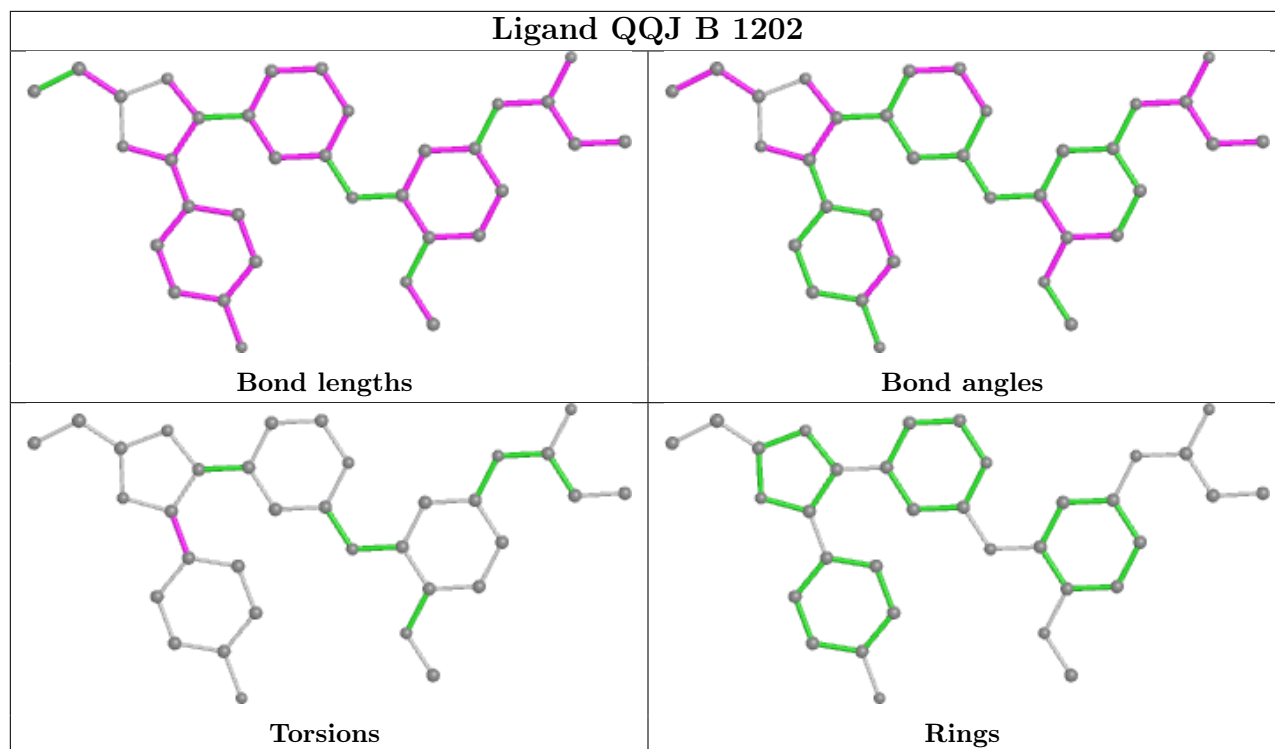
No monomer is involved in short contacts.

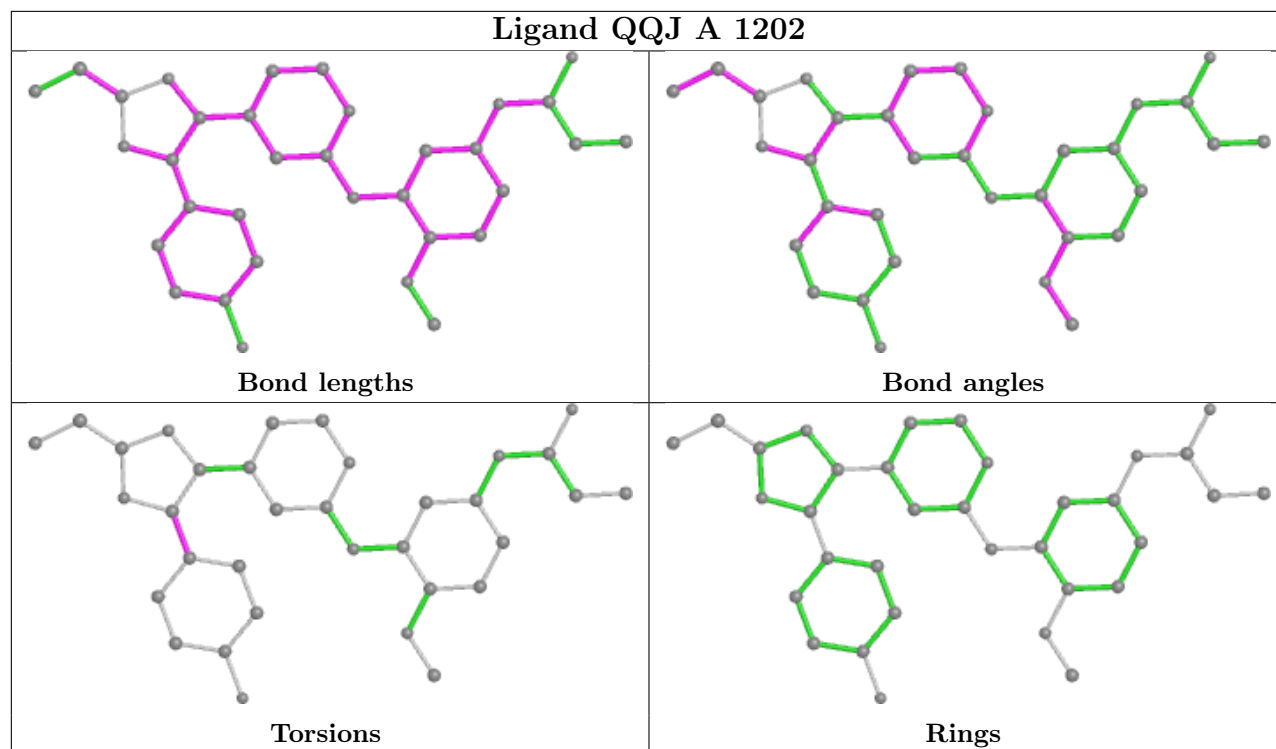
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.











4.7 Other polymers [i](#)

There are no such residues in this entry.

4.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

5 Fit of model and data

5.1 Protein, DNA and RNA chains

EDS failed to run properly - this section is therefore empty.

5.2 Non-standard residues in protein, DNA, RNA chains

EDS failed to run properly - this section is therefore empty.

5.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

5.4 Ligands

EDS failed to run properly - this section is therefore empty.

5.5 Other polymers

EDS failed to run properly - this section is therefore empty.